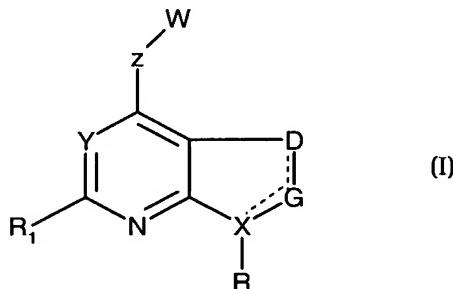


**Amendments to the claims**

This listing of claims will replace all prior versions and listings of claims in the application:

1. (Currently amended): A Compounds compound, including stereoisomers, of formula (I) including stereoisomers, prodrugs and pharmaceutically acceptable salts or solvates thereof



or a prodrug, or a pharmaceutically acceptable salt or solvate thereof, wherein  
the dashed line may represent a double bond;

R is aryl or heteroaryl, each of which may be substituted by 1 to 4 groups J selected from:  
halogen, C1-C6 alkyl, C1-C6 alkoxy, halo C1-C6 alkyl, C2-C6 alkenyl, C2-C6 alkynyl, halo C1-C6 alkoxy, -C(O)R<sub>2</sub>, nitro, hydroxy, -NR<sub>3</sub>R<sub>4</sub>, cyano or a group Z;

R<sub>1</sub> is hydrogen, C3-C7 cycloalkyl, C1-C6 alkyl, C1-C6 alkoxy, C1-C6 thioalkyl, C2-C6 alkenyl, C2-C6 alkynyl, halo C1-C6 alkyl, halo C1-C6 alkoxy, halogen, NR<sub>3</sub>R<sub>4</sub> or cyano;

R<sub>2</sub> is a C1-C4 alkyl, -OR<sub>3</sub> or -NR<sub>3</sub>R<sub>4</sub>;

R<sub>3</sub> is hydrogen or C1-C6 alkyl;

R<sub>4</sub> is hydrogen or C1-C6 alkyl;

R<sub>5</sub> is a C1-C6 alkyl, halo C1-C6 alkyl, C1-C6 alkoxy, halo C1-C6 alkoxy, C3-C7 cycloalkyl, hydroxy, halogen, nitro, cyano, -NR<sub>3</sub>R<sub>4</sub>[[;]], or -C(O)R<sub>2</sub>;

R<sub>6</sub> is a C1-C6 alkyl, halo C1-C6 alkyl, C1-C6 alkoxy, halo C1-C6 alkoxy, C3-C7 cycloalkyl, hydroxy, halogen, nitro, cyano, -NR<sub>3</sub>R<sub>4</sub>[[;]], or -C(O)R<sub>2</sub>;

R<sub>7</sub> is hydrogen, C1-C6 alkyl, halogen or halo C1-C6 alkyl;

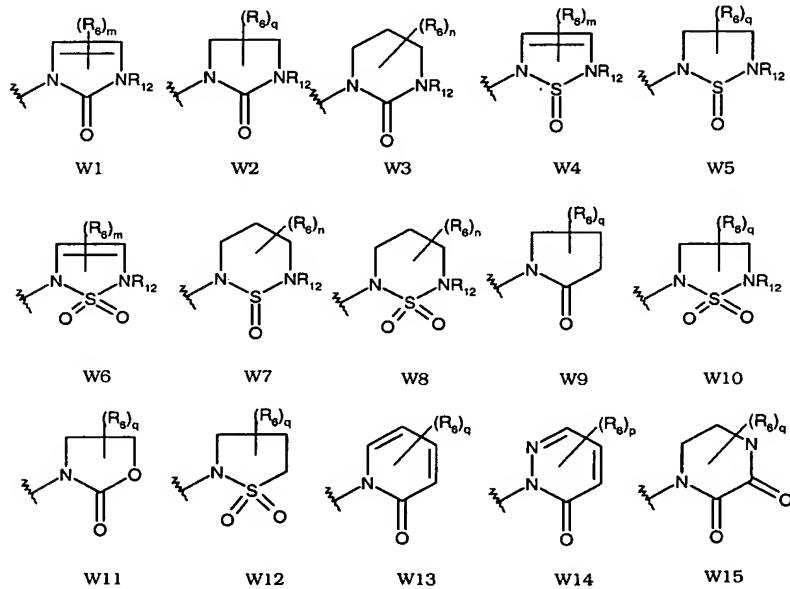
R<sub>8</sub> is hydrogen, C3-C7 cycloalkyl, C1-C6 alkyl, C2-C6 alkenyl, C2-C6 alkynyl, NR<sub>3</sub>R<sub>4</sub> or cyano;

R<sub>9</sub> is hydrogen, C3-C7 cycloalkyl, C1-C6 alkyl, C2-C6 alkenyl, C2-C6 alkynyl, NR<sub>3</sub>R<sub>4</sub> or cyano;

R<sub>10</sub> is hydrogen, C3-C7 cycloalkyl, C1-C6 alkyl, C2-C6 alkenyl, C2-C6 alkynyl, NR<sub>3</sub>R<sub>4</sub> or cyano;

$R_{11}$	is hydrogen, C3-C7 cycloalkyl, C1-C6 alkyl, C2-C6 alkenyl, C2-C6 alkynyl, $NR_3R_4$ or cyano;
$R_{12}$	is $R_3$ or $-C(O)R_2$ ;
$D$	is $CR_8R_9$ or is $CR_8$ when double bonded with $G$ ;
$G$	is $CR_{10}R_{11}$ or is $CR_{10}$ when double bonded with $D$ or is $CR_{10}$ when double bonded with $X$ when $X$ is carbon;
$X$	is carbon or nitrogen;
$Y$	is nitrogen or $-CR_7$ ;
$W$	is a 4-8 membered ring, which may be saturated or may contain one to three double bonds, and in which: <ul style="list-style-type: none"> <li>- one carbon atom is replaced by a carbonyl or <math>S(O)_m</math>; and</li> <li>- one to four carbon atoms may optionally be replaced by oxygen, nitrogen or <math>NR_{12}</math>, <math>S(O)_m</math>, carbonyl, and such ring may be further substituted by 1 to 8 <math>R_6</math> groups;</li> </ul>
$Z$	is a 5-6 membered heterocycle, which may be substituted by 1 to 8 $R_5$ groups or a phenyl ring, which may be substituted by 1 to 4 $R_5$ groups;
$m$	is an integer from 0 to 2.

2. (Currently amended): A Compounds compound according to claim 1, in which  $W$  is selected among from the following groups:

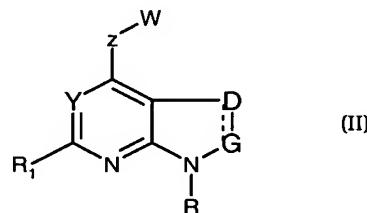


in which:

- W1 represents a 1,3-dihydro-2H-imidazol-2-one derivative;
- W2 represents a imidazolidin-2-one derivative;
- W3 represents a tetrahydropyrimidin-2(1H)-one derivative;
- W4 represents a 2,5-dihydro-1,2,5-thiadiazole 1-oxide derivative;
- W5 represents a 1,2,5-thiadiazolidine 1-oxide derivative;

W6 represents a 2,5-dihydro-1,2,5-thiadiazole 1,1-dioxide derivative;  
W7 represents a 1,2,6-thiadiazinane 1-oxide derivative;  
W8 represents a 1,2,6-thiadiazinane 1,1-dioxide derivative;  
W9 represents a pyrrolidin-2-one derivative;  
W10 represents a 2,5-dihydro-1,2,5-thiadiazolidine 1,1-dioxide derivative;  
W11 represents a 1,3-oxazolidin-2-one derivative;  
W12 represents a isothiazolidine 1,1-dioxide derivative;  
W13 represents a 2(1H)-pyridinone derivative;  
W14 represents a 3(2H)-pyridazinone;  
W15 represents a 2,3-piperazinedione derivative;  
and  
q is an integer from 0 to 4[[],];  
n is an integer from 0 to 6[[],];  
p is an integer from 0 to 3;  
m, R<sub>6</sub> and R<sub>12</sub> are defined as in claim 1; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

3. (Currently amended): A Compounds compound according to claim 1, having formula (II)



in which X is nitrogen or carbon and R, R<sub>1</sub>, Y, Z, W, D, and G have the meanings as defined in claim 1; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

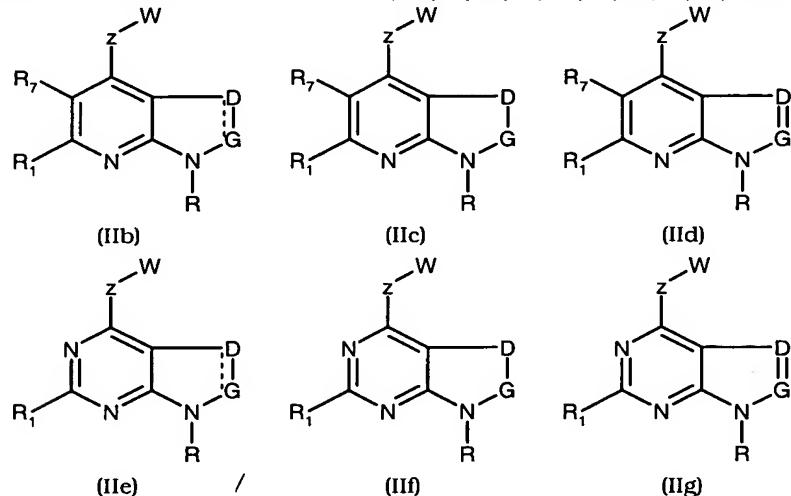
4. (Currently amended): A Compounds compound according to claim 3, of formula (II), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13, and W14; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

5. (Currently amended): A Compounds compound according to claim 3 of formula (II), in which Z is selected in the following group from: pyrimidine, pyridine, thiazole, pyrazole, triazole and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

6. (Currently amended): A Compounds compound according to any of claims from 2 to claim 3 of formula (II), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13 and W14 and in which Z is selected from the following heterocyclic groups:

pyrimidine, pyridine, thiazole, pyrazole, triazole and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

7. (Currently amended): A Compounds compound according to any one from claim 1 to claim 6 of formula (IIb), (IIc), (IId), (IIe), (IIf), and or (IIg)



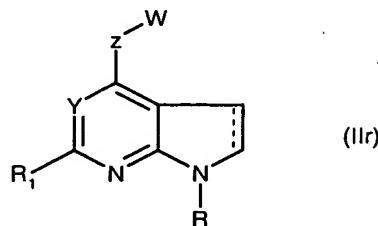
where R, R<sub>1</sub>, R<sub>7</sub>, Z, W, D, and G have the meanings as defined in claim 1; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

8. (Currently amended): A Compounds compound according to claim 7 of formula (IIb), (IIc), (IId), (IIe), (IIf) and or (IIg), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13 and W14; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

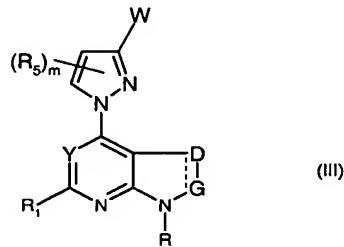
9. (Currently amended): A Compounds compound according to claims claim 7 and 8 of formula (IIb), (IIc), (IId), (IIe), (IIf) and or (IIg), in which Z is selected in the group consisting from: pyrimidine, pyridine, thiazole, pyrazole, triazole and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

10. (Currently amended): A Compounds compound according to any of claims from 7 to 9 claim 7 of formula (IIb), (IIc), (IId), (IIe), (IIf) and or (IIg), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13 and W14 and in which Z is a derivative of the following heterocyclic groups: pyrimidine, pyridine, thiazol, pyrazol, triazol and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

11. (Currently amended): A Compounds compound according to claim 7 of formula (IIr), which corresponds to the compounds a compound of formula (II), where D and G are -CH<sub>2</sub>-; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

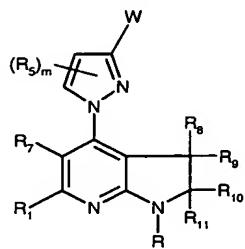


12. (Currently amended): A Compounds compound according to claim 11 of formula (IIr), in which W is selected ~~in the group consisting~~ from: W1, W2, W3, W9, W10, W11, W12, W13 and W14; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
13. (Currently amended): A Compounds compound according to ~~claims 11 and 12~~ claim 11 of formula (IIr), in which Z is selected ~~in the group consisting~~ from: pyrimidine, pyridine, thiazol, pyrazol, triazol and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
14. (Currently amended): A Compounds compound according to ~~any of claims from 11 to 13~~ claim 11 of formula (IIr), in which W is selected ~~in the group consisting~~ from: W1, W2, W3, W9, W10, W11, W12, W13 and W14 and in which Z is selected ~~in the group consisting~~ from: pyrimidine, pyridine, thiazol, pyrazol, triazol and phenyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
15. (Currently amended): A Compounds compound according to claim 3 of formula (III),

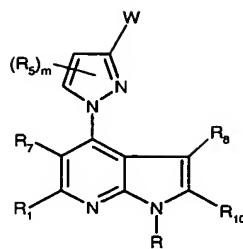


in which Z is a pyrazolyl derivative and R, R<sub>1</sub>, R<sub>5</sub>, Y, W, D, m and G have the meanings as defined in claim 1 and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

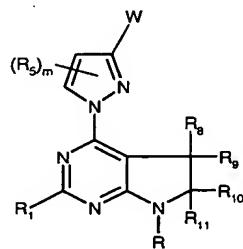
16. (Currently amended): A Compounds compound according to claim 15 of formula (IIIa), (IIIb), (IIIc) and or (IIId),



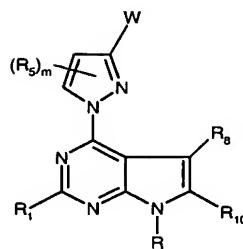
(IIIa)



(IIIb)



(IIIc)

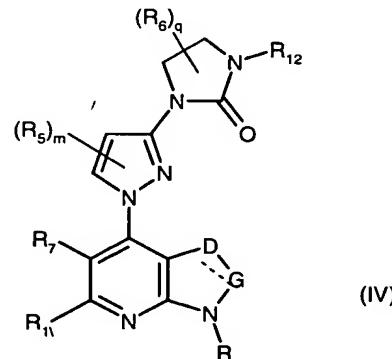


(IIId)

in which R, R<sub>1</sub>, R<sub>5</sub>, R<sub>7</sub>, R<sub>8</sub>, R<sub>9</sub>, R<sub>10</sub>, R<sub>11</sub>, W, D, m and G have the meanings as defined in claim 1; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

17. (Currently amended): A Compounds compound according to claim 16 of formula (IIIa), (IIIb), (IIIc) and or (IIId), in which W is selected in the group consisting from: W1, W2, W3, W9, W10, W11, W12, W13, W14 and R, R<sub>1</sub>, R<sub>5</sub>, R<sub>7</sub>, R<sub>8</sub>, R<sub>9</sub>, R<sub>10</sub>, R<sub>11</sub>, and m have the meanings as defined in claim 1; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

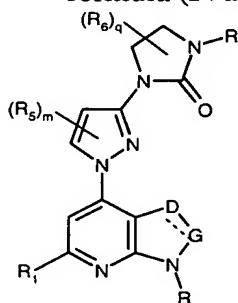
18. (Currently amended): A Compounds compound according to claim 15 of formula (IV),



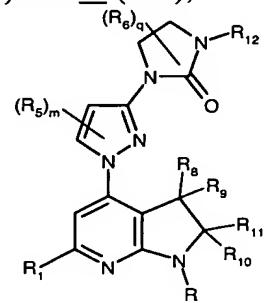
(IV)

in which R, R<sub>1</sub>, R<sub>5</sub>, R<sub>6</sub>, R<sub>7</sub>, R<sub>12</sub>, m, q, D and G have the meanings as defined in claims 1 and 2 and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

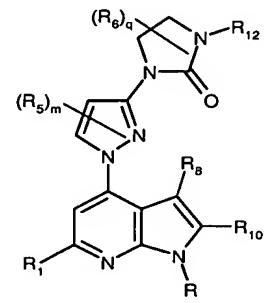
19. (Currently amended): A Compounds compound according to claim 18 of formula (IVa), (IVb) and or (IVc),



(IVa)



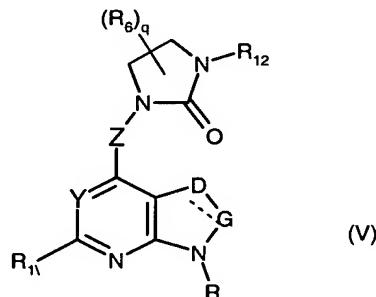
(IVb)



(IVc)

in which R, R<sub>1</sub>, R<sub>5</sub>, R<sub>6</sub>, R<sub>7</sub>, R<sub>12</sub>, m, q, D and G have the meanings as defined in claim 1 and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

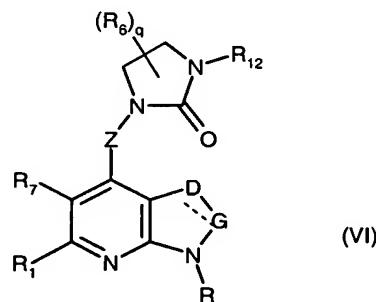
20. (Currently amended): A Compounds compound according to claim 3 of formula (V),



(V)

in which Z, R, R<sub>1</sub>, R<sub>6</sub>, q, Y, W, D and G have the meanings as defined in claims 1 and 2, and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

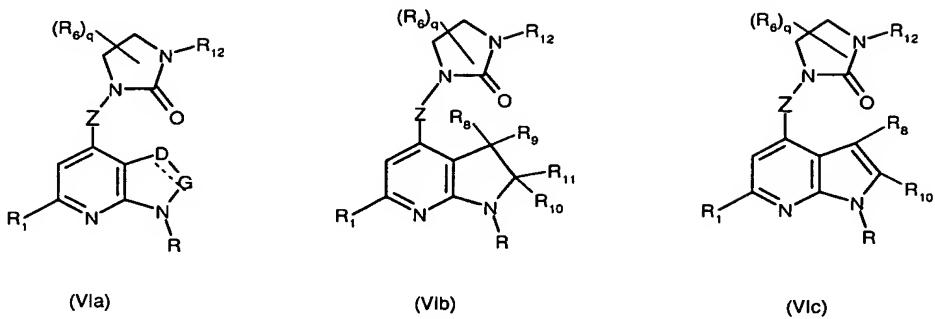
21. (Currently amended): A Compounds compound according to claim 20 of formula (VI),



(VI)

in which Z, R, R<sub>1</sub>, R<sub>6</sub>, R<sub>7</sub>, q, Y, W, D and G have the meanings as defined in claims 1 and 2, and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

22. (Currently amended): A Compounds compound according to claim 21 of formula (VIa), (VIb) and or (VIc),



in which R, R<sub>1</sub>, R<sub>6</sub>, R<sub>8</sub>, R<sub>9</sub>, R<sub>10</sub>, R<sub>11</sub>, R<sub>12</sub>, q, D and G have the meanings as defined in claims 1 and 2 and the dashed line may represent a double bond; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

23. (Currently amended): A Compounds compound according to claim 22 of formula (VIa), (VIb) and or (VIc), in which Z is selected in the group consisting from: pyrimidine, pyridine, thiazol, pyrazol, triazol and phenyl and R, R<sub>1</sub>, R<sub>6</sub>, R<sub>8</sub>, R<sub>9</sub>, R<sub>10</sub>, R<sub>11</sub>, R<sub>12</sub>, q, D and G have the meanings as defined in claim 1 and 2; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

24. (Currently amended): A Compounds compound according to any of claims from 1 to 23 of formula (I), (IIb), (He), (Id), (He), (If), (Hg), (IIr), (III), (IIIa), (IIIb), (IIIc), (IIId), (IV), (IVa), (IVb), (IVc), (V), (VI), (VIa), (VIb), (VIc) claim 1, wherein:

R<sub>1</sub> is C1-C3 alkyl group or halo C1-C3 alkyl group,

R<sub>7</sub> is hydrogen;

R<sub>8</sub>, (R<sub>9</sub>), R<sub>10</sub>, (R<sub>11</sub>) are hydrogen;

R is an aryl group selected from: 2,4-dichlorophenyl, 2-chloro-4-methylphenyl, 2-chloro-4-trifluoromethylphenyl, 2-chloro-4-methoxyphenyl, 2,4,5-trimethylphenyl, 2,4-dimethylphenyl, 2-methyl-4-methoxyphenyl, 2-methyl-4-ethoxyphenyl, 2-methyl-4-isopropoxyphenyl, 2-methyl-4-hydroxyphenyl, 2-methyl-4-chlorophenyl, 2-methyl-4-trifluoromethylphenyl, 2,4-dimethoxyphenyl, 2-methoxy-4-chlorophenyl, 3-methoxy-4-chlorophenyl, 2,5-dimethoxy-4-chlorophenyl, 2-methoxy-4-isopropylphenyl, 2-methoxy-4-trifluoromethylphenyl, 2-methoxy-4-isopropylphenyl, 2-methoxy-4-methylphenyl, 2-trifluoromethyl-4-chlorophenyl, 2,4-bis-trifluoromethylphenyl, 2-trifluoromethyl-4-methylphenyl, 2-trifluoromethyl-4-methoxyphenyl, 2-difluoromethyl-4-methoxyphenyl, 2-bromo-4-isopropylphenyl, 2-methyl-4-cyanophenyl, 2-chloro-4-cyanophenyl, 2-trifluoromethyl-4-cyanophenyl, 2-trifluoromethoxy-4-cyanophenyl, 2-ethyl-4-cyanophenyl, 2-methyl-4-trifluoromethoxyphenyl, 4-methyl-6-dimethylaminopyridin-3-yl, 2,6-bismethoxy-pyridin-3-yl, 2-methyl-6-methoxy-pyridin-3-yl, 2-trifluoromethyl-6-methoxy-pyridin-3-yl 3-chloro-5-trichloromethyl-

pyridin-2-yl, 2-methyl-4-(pyrazol-1-yl)-phenyl, 2-methoxy-4-(pyrazol-1-yl)-phenyl, 2,4,6-trimethoxyphenyl, 2-methyl-4,5-benzodioxolyl, and 2-methyl-3,4-benzodioxolyl; or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

25. (Currently amended): A Compounds compound of formula (I), (IIb), (IIc), (IId), (IIe), (IIf), (IIg), (III), (IIIA), (IIIB), (IIIC), (IIId), (IV), (IVa), (IVb), (IVc), (V), (VI), (VIa), (VIb), (VIc), according to any of claims from 1 to 24 claim 1 selected in the group consisting from:

1-{1-[1-(4-Methoxy-2-methylphenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}imidazolidin-2-one (compound 1-1);

1-{1-[1-(4-Methoxy-2-methylphenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}-3-methylimidazolidin-2-one (compound 1-2);

1-{1-[1-(2,4-Dichlorophenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}imidazolidin-2-one (compound 1-3);

1-(1-{1-[2,4-Bis(trifluoromethyl)phenyl]-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-4);

1-{1-[1-(4-Hydroxy-2-methylphenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}-2-imidazolidinone (compound 1-5);

1-Acetyl-3-(1-{6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-5);

1-Acetyl-3-(1-{6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-6);

1-(1-{1-[4-(Ethyoxy)-2-methylphenyl]-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-7);

1-[1-(6-Methyl-1-{2-methyl-4-[(1-methylethyl)oxy]phenyl}-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl)-1H-pyrazol-3-yl]-2-imidazolidinone (compound 1-8);

1-[1-(6-Methyl-1-{2-methyl-4-[(trifluoromethyl)oxy]phenyl}-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl)-1H-pyrazol-3-yl]-2-imidazolidinone (compound 1-9);

3-Methyl-4-{6-methyl-4-[3-(2-oxo-1-imidazolidinyl)-1H-pyrazol-1-yl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-1-yl}benzonitrile (compound 1-10);

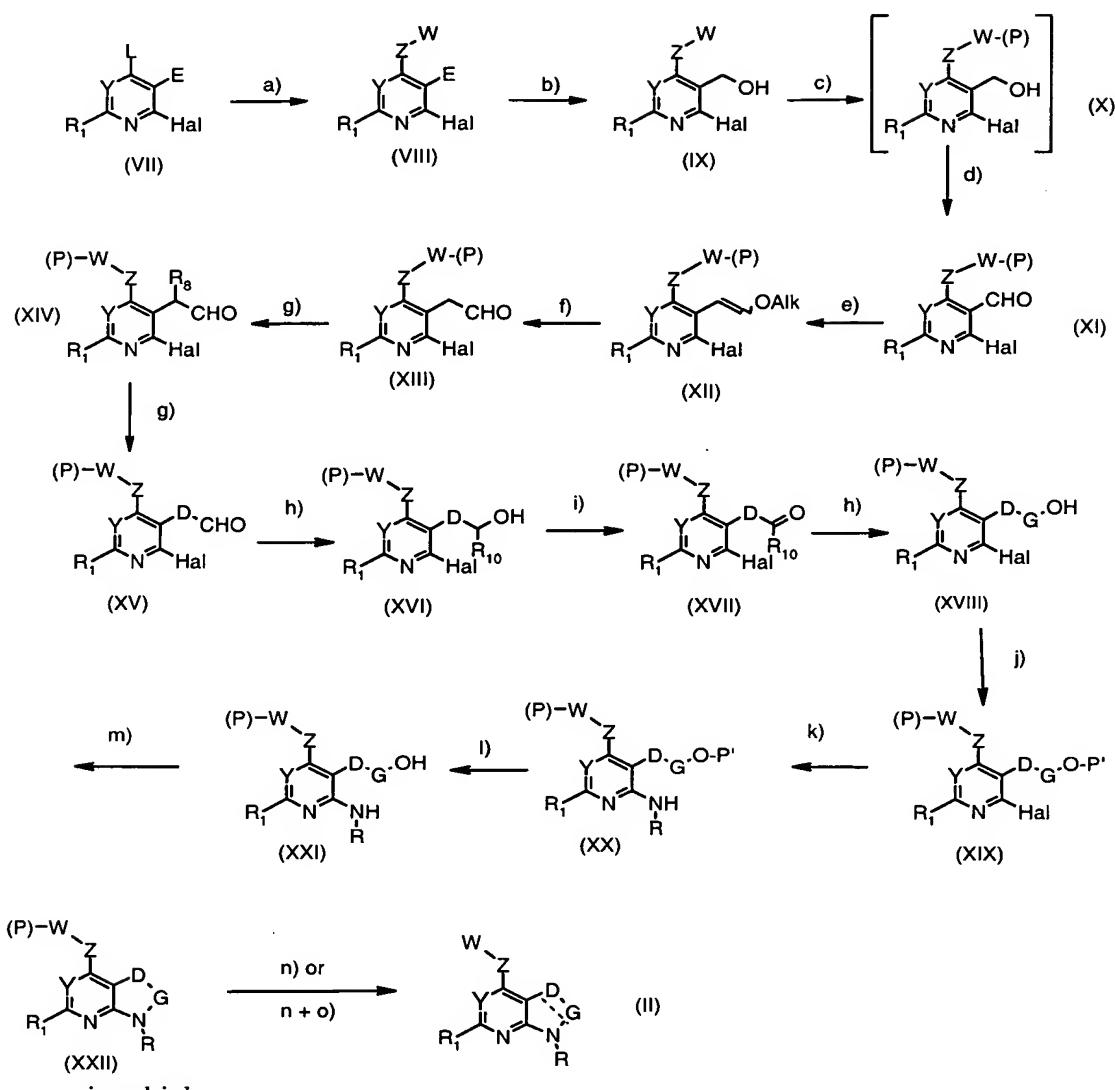
1-(1-{6-Methyl-1-[2-methyl-4-(1H-pyrazol-1-yl)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-11);

4-{6-Methyl-4-[3-(2-oxo-1-imidazolidinyl)-1H-pyrazol-1-yl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-1-yl}-3-(trifluoromethyl)benzonitrile  
(compound 1-12);  
1-(1-{1-[2-(Difluoromethyl)-4-(methyloxy)phenyl]-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone  
(compound 1-13);  
4-{6-Methyl-4-[3-(2-oxo-1-imidazolidinyl)-1H-pyrazol-1-yl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-1-yl}-3-[(trifluoromethyl)oxy]benzonitrile  
(compound 1-14);  
3-Ethyl-4-{6-methyl-4-[3-(2-oxo-1-imidazolidinyl)-1H-pyrazol-1-yl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-1-yl}benzonitrile (compound 1-15);  
1-(1-{6-Methyl-1-[2-(methyloxy)-4-(1H-pyrazol-1-yl)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-16);  
1-{1-[6-Methyl-1-(6-methyl-1,3-benzodioxol-5-yl)-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}-2-imidazolidinone (compound 1-17);  
1-(1-{6-Methyl-1-[2,4,6-tris(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-18);  
1-{1-[6-Methyl-1-(6-methyl-1,3-benzodioxol-5-yl)-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl]-1H-pyrazol-3-yl}-2-imidazolidinone (compound 1-19);  
1-(6-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-2-pyridinyl)-2-imidazolidinone (compound 1-20);  
1-(4-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-2-pyrimidinyl)-2-imidazolidinone (compound 1-21);  
1-(2-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-4-pyrimidinyl)-2-imidazolidinone (compound 1-22);  
1-(1-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-23);  
1-(1-{2,6-Dimethyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 1-24);  
1-(3-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}phenyl)-2-imidazolidinone (compound 1-25);  
1-(5-Methyl-1-{6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-b]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound

1-26);  
1-[1-(1-{4-[(difluoromethyl)oxy]-2-methylphenyl}-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridin-4-yl)-1H-pyrazol-3-yl]-2-imidazolidinone (compound 1-27);  
1-{1-[1-(4-Methoxy-2-methylphenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridin-4-yl]-1H-pyrazol-3-yl}pyrrolidin-2-one (compound 2-1);  
1-{1-[1-(4-Methoxy-2-methylphenyl)-6-methyl-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridin-4-yl]-1H-pyrazol-3-yl}tetrahydropyrimidin-2(1H)-one (compound 3-1);  
3-(1-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridin-4-yl}-1H-pyrazol-3-yl)-1,3-oxazolidin-2-one (compound 4-1);  
Methyl 5-(1-{6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridin-4-yl}-1H-pyrazol-3-yl)-1,2,5-thiadiazolidine-2-carboxylate 1,1-dioxide) (compound 5-1);  
4-[3-(1,1-Dioxido-1,2,5-thiadiazolidin-2-yl)-1H-pyrazol-1-yl]-6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridine (compound 5-2).  
4-[3-(1,1-Dioxido-2-isothiazolidinyl)-1H-pyrazol-1-yl]-6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridine (compound 6-1);  
3-Methyl-1-(1-{6-methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridin-4-yl}-1H-pyrazol-3-yl)-2(1H)-pyridinone (compound 7-1);  
2-(1-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridin-4-yl}-1H-pyrazol-3-yl)-3(2H)-pyridazinone (compound 8-1);  
1-(1-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridin-4-yl}-1H-pyrazol-3-yl)-1,3-dihydro-2H-imidazol-2-one (compound 9-1);  
1-(1-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-1H-pyrrolo[2,3-*b*]pyridin-4-yl}-1H-pyrazol-3-yl)-2-imidazolidinone (compound 10-1);  
1-(6-{6-Methyl-1-[2-methyl-4-(methyloxy)phenyl]-2,3-dihydro-1H-pyrrolo[2,3-*b*]pyridin-4-yl}-3-pyridinyl)-2-imidazolidinone (compound 11-1); and  
1-{1-[7-(2,4-Dichlorophenyl)-2-methyl-6,7-dihydro-5H-pyrrolo[2,3-*d*]pyrimidin-4-yl]-1H-pyrazol-3-yl}-2-pyrrolidinone (compound 11-2); or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.

26. (Currently amended): A process for the preparation of a compound the compounds of formula (II), starting from a compound compounds of formula (VII), comprising the following steps as in Scheme 1:

Scheme 1



- step a stands for conversion of the leaving group L, selected in a group consisting from: halogen or reactive residue of sulphonic acid (e.g. mesylate, tosylate), preferably chloride, in the compounds (VIII), by reaction with the suitable Z-W derivative;
- step b stands for reduction of the ester group (E) with a suitable reducing agent (such as ~~DIBAL-H~~) to hydroxy group of compounds (IX);
- step c stands for suitable protection of an NH group eventually present in W group with a P group, such as a p-methoxybenzyl group;
- step d stands for oxidation of the hydroxy group with a suitable oxidizing agent (such as ~~Dess-Martin periodinane~~) to the aldehyde group of compounds (XI);

steps e + f stands for formation of the aldehyde group of compounds (XIII) by Wittig reaction in the usual conditions, through formation of enol ether followed by acid hydrolysis (step f);

step g stands for the optional alkylation of the  $\alpha$  position of the aldehyde by deprotonation with a suitable base (such as  $\text{LiN}(\text{SiMe}_3)_2$ ), followed by the addition of a suitable alkylating agent (such as  $\text{MeI}$ ) to form the alkylated aldehyde of compounds (XIV), (XV);

step h stands for the conversion of the aldehyde group group by a Grignard reagent (such as  $\text{MeMgBr}$ ) into an alcohol group of compounds (XVI) and (XVIII);

step i stands for oxidation of the hydroxy group with a suitable oxidizing agent (such as Dess-Martin periodinane) to the ketone group of compounds (XVII);

step j stands for conversion of the hydroxy group in the suitable protecting group of compounds (XIX) (such as TBS:tert-butyldimethylsilyl);

step k stands for a Buchwald coupling reaction with the suitable amine  $\text{RNH}_2$  to give the compounds of formula (XX);

step l stands for the deprotection reaction to give the hydroxy group of compounds (XXI);

step m stands for intramolecular cyclisation after conversion of the hydroxy group of compounds (XXI) in a suitable leaving group (such as bromide, by reaction with  $\text{CBr}_4$  and  $\text{PPh}_3$ ) to give the cyclized compounds (XXII);

step n stands for the deprotection reaction of the protected NH group eventually present in W group, to give final compounds (II); and

step o stands for oxidation by a suitable oxidizing agent (such as DDQ) in order to give formation of the double bond of compounds (II), when D is  $\text{CHR}_8$  and G is  $\text{CHR}_{10}$ .

27-32. (Cancelled)

33. (Currently amended): A pharmaceutical composition comprising a compound according to any of claims from 1 to 25 claim 1, or a prodrug, or a pharmaceutically acceptable salt or solvate thereof, in admixture with one or more physiologically acceptable carriers or excipients.

34. (Currently amended): A method for the treatment of a mammal, including man, in particular in the treatment of a condition conditions mediated by CRF (corticotropin-releasing factor), comprising administration of an effective amount of a compound according to any of claims from 1 to 25

claim 1, or a prodrug, or a pharmaceutically acceptable salt or solvate thereof, to a mammal in need thereof.

35. (Currently amended): A method, according to claim 34, in the treatment of depression and anxiety, comprising administration of an effective amount of a compound according to ~~any of claims 1 to 25~~claim 1, or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.
36. (Currently amended): A method, according to claim 34, in the treatment of IBS (irritable bowel disease) and IBD (inflammatory bowel disease), comprising administration of an effective amount of a compound according to ~~any of claims 1 to 25~~claim 1, or a prodrug, or a pharmaceutically acceptable salt or solvate thereof.